

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041520\
 Data File : BG044988.D
 Acq On : 15 Apr 2020 18:57
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 16 05:01:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 17:57:00 2020
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	97	0.00
2	1,4-Dioxane	40.000	43.276	-8.2	105	0.00
3	Pyridine	40.000	40.849	-2.1	103	0.00
4	n-Nitrosodimethylamine	40.000	41.510	-3.8	105	0.00
5 S	2-Fluorophenol	80.000	82.076	-2.6	103	0.00
6	Aniline	40.000	39.267	1.8	96	0.00
7 S	Phenol-d6	80.000	80.209	-0.3	98	0.00
8	2-Chlorophenol	40.000	39.782	0.5	98	0.00
9	Benzaldehyde	40.000	42.479	-6.2	101	0.00
10 C	Phenol	40.000	40.302	-0.8	99	0.00
11	bis(2-Chloroethyl)ether	40.000	39.725	0.7	98	0.00
12	1,3-Dichlorobenzene	40.000	40.153	-0.4	100	0.00
13 C	1,4-Dichlorobenzene	40.000	39.527	1.2	98	0.00
14	1,2-Dichlorobenzene	40.000	40.149	-0.4	98	0.00
15	Benzyl Alcohol	40.000	39.034	2.4	96	0.00
16	2,2'-oxybis(1-Chloropropane)	40.000	39.946	0.1	98	0.00
17	2-Methylphenol	40.000	39.445	1.4	98	0.00
18	Hexachloroethane	40.000	38.536	3.7	94	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	40.084	-0.2	98	0.00
20	3+4-Methylphenols	40.000	38.886	2.8	95	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	92	0.00
22	Acetophenone	40.000	40.108	-0.3	96	0.00
23 S	Nitrobenzene-d5	80.000	80.943	-1.2	98	0.00
24	Nitrobenzene	40.000	40.285	-0.7	96	0.00
25	Isophorone	40.000	40.302	-0.8	94	0.00
26 C	2-Nitrophenol	40.000	41.905	-4.8	100	0.00
27	2,4-Dimethylphenol	40.000	39.043	2.4	93	0.00
28	bis(2-Chloroethoxy)methane	40.000	40.510	-1.3	95	0.00
29 C	2,4-Dichlorophenol	40.000	40.521	-1.3	96	0.00
30	1,2,4-Trichlorobenzene	40.000	39.861	0.3	95	0.00
31	Naphthalene	40.000	40.197	-0.5	94	0.00
32	Benzoic acid	40.000	38.504	3.7	96	0.00
33	4-Chloroaniline	40.000	39.658	0.9	95	0.00
34 C	Hexachlorobutadiene	40.000	40.143	-0.4	96	0.00
35	Caprolactam	40.000	38.762	3.1	91	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.211	2.0	94	0.00
37	2-Methylnaphthalene	40.000	40.094	-0.2	94	0.00
38	1-Methylnaphthalene	40.000	39.874	0.3	94	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	91	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	41.672	-4.2	97	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.434	-3.6	94	0.00
42 S	2,4,6-Tribromophenol	80.000	81.840	-2.3	96	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.802	-2.0	94	0.00
44	2,4,5-Trichlorophenol	40.000	40.122	-0.3	93	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041520\
 Data File : BG044988.D
 Acq On : 15 Apr 2020 18:57
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 16 05:01:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 17:57:00 2020
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45 S	2-Fluorobiphenyl	80.000	82.779	-3.5	95	0.00
46	1,1'-Biphenyl	40.000	41.065	-2.7	95	0.00
47	2-Chloronaphthalene	40.000	39.925	0.2	94	0.00
48	2-Nitroaniline	40.000	41.197	-3.0	98	0.00
49	Acenaphthylene	40.000	40.585	-1.5	94	0.00
50	Dimethylphthalate	40.000	40.053	-0.1	93	0.00
51	2,6-Dinitrotoluene	40.000	40.298	-0.7	95	0.00
52 C	Acenaphthene	40.000	40.700	-1.8	95	0.00
53	3-Nitroaniline	40.000	39.855	0.4	93	0.00
54 P	2,4-Dinitrophenol	40.000	41.672	-4.2	100	0.00
55	Dibenzofuran	40.000	39.975	0.1	93	0.00
56 P	4-Nitrophenol	40.000	40.506	-1.3	95	0.00
57	2,4-Dinitrotoluene	40.000	39.943	0.1	95	0.00
58	Fluorene	40.000	40.065	-0.2	93	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.846	-2.1	96	0.00
60	Diethylphthalate	40.000	39.343	1.6	92	0.00
61	4-Chlorophenyl-phenylether	40.000	39.945	0.1	94	0.00
62	4-Nitroaniline	40.000	40.092	-0.2	94	0.00
63	Azobenzene	40.000	40.409	-1.0	95	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	91	0.00
65	4,6-Dinitro-2-methylphenol	40.000	40.780	-2.0	98	0.00
66 c	n-Nitrosodiphenylamine	40.000	40.071	-0.2	93	0.00
67	4-Bromophenyl-phenylether	40.000	40.838	-2.1	95	0.00
68	Hexachlorobenzene	40.000	39.822	0.4	94	0.00
69	Atrazine	40.000	41.757	-4.4	94	0.00
70 C	Pentachlorophenol	40.000	41.280	-3.2	94	0.00
71	Phenanthrene	40.000	40.812	-2.0	95	0.00
72	Anthracene	40.000	40.585	-1.5	95	0.00
73	Carbazole	40.000	39.315	1.7	95	0.00
74	Di-n-butylphthalate	40.000	42.259	-5.6	94	0.00
75 C	Fluoranthene	40.000	42.739	-6.8	96	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	96	0.00
77	Benzidine	40.000	45.677	-14.2	103	0.00
78	Pyrene	40.000	38.904	2.7	97	0.00
79 S	Terphenyl-d14	80.000	83.524	-4.4	98	0.00
80	Butylbenzylphthalate	40.000	40.551	-1.4	99	0.00
81	Benzo(a)anthracene	40.000	41.071	-2.7	101	0.00
82	3,3'-Dichlorobenzidine	40.000	41.082	-2.7	100	0.00
83	Chrysene	40.000	40.615	-1.5	100	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.767	-1.9	100	0.00
85 c	Di-n-octyl phthalate	40.000	41.491	-3.7	101	0.00
86	Indeno(1,2,3-cd)pyrene	40.000	41.376	-3.4	103	0.00
87 I	Perylene-d12	20.000	20.000	0.0	98	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041520\
Data File : BG044988.D
Acq On : 15 Apr 2020 18:57
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Apr 16 05:01:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 15 17:57:00 2020
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	41.301	-3.3	105	0.00
89	Benzo(k)fluoranthene	40.000	40.463	-1.2	101	0.00
90 C	Benzo(a)pyrene	40.000	40.556	-1.4	102	0.00
91	Dibenzo(a,h)anthracene	40.000	40.426	-1.1	103	0.00
92	Benzo(q,h,i)perylene	40.000	40.613	-1.5	104	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0